

each R₁ is independently selected from the group consisting of H; halogen; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; --COR₄ where R₄ is H, C₁₋₄ alkyl or C₁₋₄ alkoxy; C₃₋₆ cycloalkyl; aryl; heteroaryl; cyano; nitro; trihalomethyl; oxo; or -(CH₂)_n-X-(CH₂)_m-(R₅)_o where X is O, S or N, n is 0-3, m is 0-3, o is 0-1, and R₅ is methyl or H₁₋₂;

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each R₂ and each R₃ are independently selected from the group consisting of H; halogen; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; --COR₄ where R₄ is H; C₁₋₄ alkyl or C₁₋₄ alkoxy; C₃₋₆ cycloalkyl; aryl; heteroaryl; cyano; nitro; trihalomethyl; oxo; or -(CH₂)_n-X-(CH₂)_m-(R₅)_o where X is O, S or N, n is 0-3, m is 0-3, o is 0-1, and R₅ is methyl or H₁₋₂; or an R₂ and an R₃ together comprise a saturated, partly saturated, or unsaturated ring structure having the formula -(C(R₆)_p)_q-X_s-(C(R₆)_p)_r-X_t-(C(R₆)_p)_u where each R₆ is independently selected from the group consisting of H; halogen; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; --COR₄ where R₄ is H, C₁₋₄ alkyl or C₁₋₄ alkoxy; C₃₋₆ cycloalkyl; aryl; heteroaryl; cyano; nitro; trihalomethyl and oxo where each p is independently 1 or 2, q is 0-5, r is 0-5, u is 0-5; each X is independently O, S, or N and s is 0 or 1; provided that q + r + u + s + t is less than 6;

Y is selected from the group consisting of O; S; N; --(C(R₇)_z)_s—where each R₇ is independently as previously defined for R1, each z is independently 1-2, and s is 1-3; --CH=; --CH=CH--; or Y₁CH₂—where Y₁ is O, N, or S; and the dotted lines are optional double bonds, with the proviso that if the ring including Y is a cyclohexane ring or a heterocyclic 5 member ring said ring is not fully unsaturated, and that if Y is O, N or S, the ring including Y contains at least one said double bond,

said compound further having selective agonist activity at the α2B or α2B/α2C adrenergic receptor subtype(s) over the α2A adrenergic receptor subtype, and all pharmacologically acceptable salts, esters, stereoisomers and racemic mixtures thereof.

2. (Amended) The compound of claim 1 in which the ring including Y has either a single double bond or no double bond, except that when an R₂ and an R₃ together comprise a saturated, unsaturated or partly saturated ring structure said Y-including ring optionally shares an additional double bond with said condensed ring, provided Y is not S, O, or N.

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9. (Amended) The compound of claim 2, in which each R₂ and each R₃ are independently selected from the group consisting of: H; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; halide; trihalomethyl; cycloalkyl; (CH₂)_n-X-(CH₂)_m-(R₅)_o, where X is O, S or N, n is 0-3, m is 0-3, o is 0-1, and R₅ is methyl or H₁₋₂; or an R₂ and an R₃ together comprise a saturated, partly saturated, or unsaturated ring structure having the formula -(C(R₆)_p)_q-X_s-(C(R₆)_p)_r-X_t-(C(R₆)_p)_u where each R₆ is independently selected from the group consisting of H; halogen; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; --COR₄ where R₄ is H, C₁₋₄ alkyl or C₁₋₄ alkoxy; C₃₋₆ cycloalkyl; aryl; heteroaryl; cyano; nitro; trihalomethyl; and oxo where each p is independently 1 or 2, q is 0-4, r is 0-4, u is 0-4; each X is independently O, S, or N, s is 0 or 1, and q + s + r + t + u = 3 or 4.

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10. (Amended) The compound of claim 3, in which each R₂ and each R₃ are independently selected from the group consisting of: H; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; halide; trihalomethyl; cycloalkyl; (CH₂)_n-X-(CH₂)_m-(R₅)_o, where X is O, S or N, n is 0-3, m is 0-3, o is 0-1, and R₅ is methyl or H₁₋₂; or an R₂ and an R₃ together comprise a saturated, partly saturated, or unsaturated ring structure having the formula -(C(R₆)_p)_q-X_s-(C(R₆)_p)_r-X_t-(C(R₆)_p)_u where each R₆ is independently selected from the group consisting of H; halogen; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; --COR₄ where R₄ is H, C₁₋₄ alkyl or C₁₋₄ alkoxy; C₃₋₆ cycloalkyl; aryl; heteroaryl; cyano; nitro; trihalomethyl; and oxo where each p is independently 1 or 2, q is 0-4, r is 0-4, u is 0-4; each X is independently O, S, or N, s is 0 or 1, and q + s + r + t + u = 3 or 4.

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71. (Amended) The compound of claim 53 in which an R₂ and an R₃ together comprise a saturated, partly saturated, or unsaturated ring structure having the formula -(C(R₆)_p)_q-X_s-(C(R₆)_p)_r-X_t-(C(R₆)_p)_u where each R₆ is independently selected from the group consisting of H; halogen; C₁₋₄ alkyl; C₁₋₄ alkenyl; C₁₋₄ alkynyl; --COR₄ where R₄ is H, C₁₋₄ alkyl or C₁₋₄ alkoxy; C₃₋₆ cycloalkyl; aryl; heteroaryl; cyano; nitro; trihalomethyl; and oxo where each p is independently 1 or 2, q is 0-4, r is 0-4, u is 0-4; each X is independently O, S, or N, s is 0 or 1, and q + s + r + t + u = 3 or 4.